



an enzyme entry in the MetaCyc Pathway Details page, enzymes can be added to the Function Cart for further analysis.

Clicking on a Reaction ID from the MetaCyc Pathway Details page (Figure 1(ii)) will lead to a MetaCyc Reaction Detail page as shown in Figure 2(i). Clicking on a gene count will lead to a list of genomes having genes annotated with the corresponding enzyme (Figure 2(ii)). Genomes can be added to the Genome Cart for further analysis. Clicking on the gene count to view the actual genes.

### MetaCyc Reaction Detail

[DHBDEHYD-RXN](#)

Unique ID	DHBDEHYD-RXN
EC Number	<a href="#">EC:1.3.1.28</a>
Is Orphan?	No
In Pathway	<a href="#">PWY-5901</a> : 2,3-dihydroxybenzoate biosynthesis
Definition	DIHYDRO-DIOH-BENZOATE%NAD => 2-3-DIHYDROXYBENZOATE%NADH%PROTON
RHEA	23824

### Isolate Genomes with EC:1.3.1.28

Genomes with 2,3-dihydro-2,3-dihydroxybenzoate dehydrogenase..

Domains(D): \* = Microbiome,  
B = Bacteria, A = Archaea, E = Eukarya, P = Plasmids, G = GFragment, V = Viruses.  
Genome Completion(C): F = Finished, P = Permanent Draft, D = Draft.

[Add Selected to Genome Cart](#)
[Select All](#)
[Clear All](#)

Filter column: Domain Filter text:  [Apply](#)

[Export](#)
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[Column Selector](#)
[Select Page](#)
[Deselect Page](#)

Select	Domain	Status	Genome	Gene Count
<input type="checkbox"/>	A	D	<a href="#">Candidatus Pacearchaeota archaeon RBG_13_36_9</a>	<a href="#">1</a>
<input type="checkbox"/>	A	D	<a href="#">putative Thaumarchaeota genus Nitrosopumilus sp. v2 (bin U_52180)</a>	<a href="#">1</a>
<input type="checkbox"/>	B	F	<a href="#">Acinetobacter baumannii R612</a>	<a href="#">1</a>
<input type="checkbox"/>	B	P	<a href="#">Brucella melitensis bv. 1 M111</a>	<a href="#">1</a>
<input type="checkbox"/>	B	P	<a href="#">Escherichia coli ARS4 2123</a>	<a href="#">1</a>

**Figure 2.** MetaCyc Reaction and Associated Genomes and Genes

Near the top of the MetaCyc Pathways page (Figure 1(i)) there is a "MetaCyc Compounds" link leads to a list of all MetaCyc compounds loaded in IMG (Figure 3(ii)). Click on a Compound ID to view the detailed definition of the compound including compound name, molecular weight and SMILES as shown in Figure 3(ii).

**MetaCyc Compound List** (i)

Filter column:  Filter

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Column Selector

Compound ID	Compound Name	CHEBI	KEGG LIGAND	Formula	Mol. Weight
<a href="#">CPD0-1659</a>	$\alpha$ -allenylagmatine	-	-		168.241
<a href="#">CPD0-1680</a>	$\alpha$ -ethynylagmatine	-	-		154.214
<a href="#">ALPHAALPHA-TREHALOSE-6,6-BISMYCOLATE</a>	$\alpha$ , $\alpha'$ -trehalose 6,6'-bismycolate	-	-		2582.383
<a href="#">TREHALOSE</a>	$\alpha$ , $\alpha$ -trehalose	-	-		342.299
<a href="#">TREHALOSE-6P</a>	$\alpha$ , $\alpha$ -trehalose 6-phosphate	-	-		420.263
<a href="#">ALPHABETA-DIDEHYDROTRYPTOPHAN</a>	$\alpha$ , $\beta$ -didehydrotryptophan	<a href="#">CHEBI:15802</a>	-		201.204
<a href="#">CPD-7131</a>	$\alpha$ , $\beta$ -dihydroxanthohumol	-	-		355.41

**MetaCyc Compound Detail** (ii)

Unique ID	ALPHABETA-DIDEHYDROTRYPTOPHAN
Common Name	$\alpha$ , $\beta$ -didehydrotryptophan
Mol. Weight	201.204
SMILES	<chem>C2(=C(C=C(N)C(=O)[O-])C1(C=CC=CC=1N2))</chem>
DB Source	Metacyc
CHEBI	<a href="#">15802</a>

**Figure 3.** MetaCyc Compounds.